Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

- 1. (canceled)
- 2. (currently amended) A single source precursor according to claim 1, for the deposition of ternary chalcopyrite materials, said single source precursor having a structural formula selected from the group consisting of

$$\begin{bmatrix} & & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & &$$

wherein L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom, M' is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl, vinyl, perfluoro aryl, silane, and carbamato groups, said single source precursor excluding

 $[{P(C_6H_5)_3}_2Cu(S-C_2H_5)_2In(S-C_2H_5)_2],$

 $\underline{[\{P(C_6H_5)_3\}_2Cu(Se-C_2H_5)_2In(Se-C_2H_5)_2]},$

 $[\{P(C_6H_5)_3\}_2Cu(S(i-C_4H_9))_2In(S(i-C_4H_9))_2],$

 $[{P(C_6H_5)_3}_2Cu(Se(i-C_4H_9))_2In(Se(i-C_4H_9))_2],$

 $\underline{[\{P(C_6H_5)_3\}_2Ag(Cl)(SC\{O\}CH_3)In(SC\{O\}CH_3)_2]},$

 $\label{eq:continuous} \mbox{$[\{P(C_6H_5)_3\}_2Ag(Cl)(SC\{O\}C_6H_5)In(SC\{O\}C_6H_5)_2],$}$

 $[{P(C_6H_5)_3}_2Ag(SC{O}CH_3)_2In(SC{O}CH_3)_2],$

 $[\{P(C_6H_5)_3\}_2Ag(SC\{O\}C_6H_5)_2In(SC\{O\}C_6H_5)_2],$

 $\underline{[\{P(C_6H_5)_3\}_2Cu(SC\{O\}C_6H_5)_2In(SC\{O\}C_6H_5)_2]},$

 $[\{P(C_6H_5)_3\}_2Cu(SC\{O\}C_6H_5)_2Ga(SC\{O\}C_6H_5)_2],$

 $\underline{[\{P(C_6H_5)_3\}_2Ag(SC\{O\}C_6H_5)_2Ga(SC\{O\}C_6H_5)_2], and}$

 $\underline{[\{P(C_6H_5)_3\}_2Ag(SC\{O\}CH_3)_2Ga(SC\{O\}CH_3)_2]}.$

- 3. (original) A single source precursor according to claim 2, said singles source precursor being a liquid at room temperature.
- 4. (original) A single source precursor according to claim 3, said single source precursor being soluble in polar organic solvents and in non-polar organic solvents.
- 5. (original) A single source precursor according to claim 2, of the formula $[\{P(n-C_4H_9)_3\}_2Cu(Se-C_6H_5)_2In(Se-C_6H_5)_2].$
- 6. (original) A single source precursor according to claim 2, of the formula $[\{P(n-C_4H_9)_3\}_2Ag(S-C_2H_5)_2In(S-C_2H_5)_2].$
- 7. (original) A single source precursor according to claim 2, of the formula $[\{P(n-C_4H_9)_3\}_2Cu(S-C_2H_5)_2In(S-C_2H_5)_2].$
- 8. (original) A single source precursor according to claim 2, of the formula $[\{P(n-C_4H_9)_3\}_2Cu(S-C_3H_7)_2In(S-C_3H_7)_2].$
- 9. (original) A single source precursor according to claim 2, of the formula $[\{P(C_6H_5)_3\}_2Ag(S-CH_3)_2In(S-CH_3)_2].$
- 10. (original) A single source precursor according to claim 2, said single source precursor being effective to yield a I-III-VI₂ ternary chalcopyrite material upon heating or pyrolysis of said single source precursor at a temperature less than about 500°C.
- 11. (original) A single source precursor according to claim 2, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 1.5 eV between a conduction band and a valence band thereof.
- 12. (original) A single source precursor according to claim 11, said ternary chalcopyrite material being CuInS₂.

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- 13. (original) A single source precursor according to claim 2, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 2 eV between a conduction band and a valence band thereof.
- 14. (original) A single source precursor according to claim 13, said ternary chalcopyrite material being CuGaS₂.
- 15. (original) A single source precursor according to claim 2, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of 1.5-2 eV between a conduction band and a valence band thereof, said ternary chalcopyrite material being Cu(In:Ga)(S:Se)₂.
- 16. (currently amended) A single source precursor according to claim 1, for the deposition of ternary chalcopyrite materials, said single source precursor having a structural formula selected from the group consisting of

$$L \longrightarrow M! \longrightarrow M! \longrightarrow M! \longrightarrow R$$
 and
$$L \longrightarrow M! \longrightarrow R$$

$$R \longrightarrow R$$

wherein L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom, M' is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl, vinyl, perfluoro aryl, silane, and carbamato groups.

17. (original) A single source precursor according to claim 16, said single source precursor being effective to yield a I-III-VI₂ ternary chalcopyrite material upon heating or pyrolysis of said single source precursor at a temperature less than about 500°C.

- 18. (original) A single source precursor according to claim 16, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 1.5 eV between a conduction band and a valence band thereof.
- 19. (original) A single source precursor according to claim 18, said ternary chalcopyrite material being CuInS₂.
- 20. (original) A single source precursor according to claim 16, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 2 eV between a conduction band and a valence band thereof.
- 21. (original) A single source precursor according to claim 20, said ternary chalcopyrite material being CuGaS₂.
- 22. (original) A single source precursor according to claim 16, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of 1.5-2 eV between a conduction band and a valence band thereof, said ternary chalcopyrite material being Cu(In:Ga)(S:Se)₂.
- 23. (currently amended) A single source precursor according to claim 1, for the deposition of ternary chalcopyrite materials, said single source precursor having a structural formula selected from the group consisting of

wherein L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom, M' is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl, vinyl, perfluoro alkyl, perfluoro aryl, silane, and carbamato groups.

- 24. (original) A single source precursor according to claim 23, said single source precursor being effective to yield a I-III-VI2 ternary chalcopyrite material upon heating or pyrolysis of said single source precursor at a temperature less than about 500°C.
- 25. (original) A single source precursor according to claim 23, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 1.5 eV between a conduction band and a valence band thereof.
- 26. (original) A single source precursor according to claim 25, said ternary chalcopyrite material being CuInS₂.
- 27. (original) A single source precursor according to claim 23, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 2 eV between a conduction band and a valence band thereof.
- 28. (currently amended) A single source precursor according to claim 27, said ternary chalcopyrite material being CuGaS CuGaS2.
- 29. (original) A single source precursor according to claim 23, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of 1.5-2 eV between a conduction band and a valence band thereof, said ternary chalcopyrite material being $Cu(In:Ga)(S:Se)_2$.
- 30. (currently amended) A single source precursor according to claim 1, for the deposition of ternary chalcopyrite materials, said single source precursor having a structural formula selected from the group consisting of

$$L \longrightarrow M! \longrightarrow M! \longrightarrow R$$
 and
$$L \longrightarrow M! \longrightarrow R$$

$$R$$

$$10$$

wherein L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom, M' is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl, vinyl, perfluoro aryl, silane, and carbamato groups.

- 31. (original) A single source precursor according to claim 30, said single source precursor being effective to yield a I-III-VI₂ ternary chalcopyrite material upon heating or pyrolysis of said single source precursor at a temperature less than about 500°C.
- 32. (original) A single source precursor according to claim 30, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 1.5 eV between a conduction band and a valence band thereof.
- 33. (original) A single source precursor according to claim 32, said ternary chalcopyrite material being CuInS₂.
- 34. (original) A single source precursor according to claim 30, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of about 2-2.4 eV between a conduction band and a valence band thereof.
- 35. (original) A single source precursor according to claim 34, said ternary chalcopyrite material being CuGaS₂.
- 36. (original) A single source precursor according to claim 30, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of 1.5-2 eV between a conduction band and a valence band thereof, said ternary chalcopyrite material being Cu(In:Ga)(S:Se)₂.
- 37. (currently amended) A single source precursor-according to claim 1, having three E-R groups for the deposition of ternary chalcopyrite materials, said single source precursor having the empirical formula $[\{L\}_nM'(ER)_x(X)_y(R)_zM'']$, wherein x is 3, x+y+z=4, n is greater than or equal to 1, L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom,

M is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl, vinyl, perfluoro aryl, silane, and carbamato groups, said single source precursor excluding

 $[{P(C_6H_5)_3}_2Cu(S-C_2H_5)_2In(S-C_2H_5)_2],$

 $[{P(C_6H_5)_3}_2Cu(Se-C_2H_5)_2In(Se-C_2H_5)_2],$

 $[{P(C_6H_5)_3}_2Cu(S(i-C_4H_9))_2In(S(i-C_4H_9))_2],$

 $[{P(C_6H_5)_3}_2Cu(Se(i-C_4H_9))_2In(Se(i-C_4H_9))_2],$

 $[{P(C_6H_5)_3}_2Ag(Cl)(SC{O}CH_3)In(SC{O}CH_3)_2],$

 $[{P(C_6H_5)_3}_2Ag(Cl)(SC{O}_6H_5)In(SC{O}_6H_5)_2],$

 $[{P(C_6H_5)_3}_2Ag(SC{O}CH_3)_2In(SC{O}CH_3)_2],$

 $[{P(C_6H_5)_3}_2Ag(SC{O}C_6H_5)_2In(SC{O}C_6H_5)_2],$

 $[{P(C_6H_5)_3}_2Cu(SC{O}C_6H_5)_2In(SC{O}C_6H_5)_2],$

 $[{P(C_6H_5)_3}_2Cu(SC{O}C_6H_5)_2Ga(SC{O}C_6H_5)_2],$

 $[{P(C_6H_5)_3}_2Ag(SC{O}_6H_5)_2Ga(SC{O}_6H_5)_2], and$

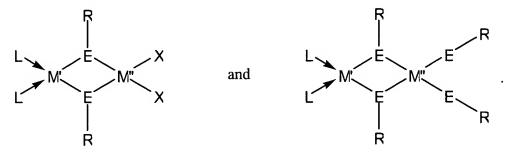
 $[{P(C_6H_5)_3}_2Ag(SC{O}CH_3)_2Ga(SC{O}CH_3)_2].$

- 38. (original) A single source precursor for the deposition of ternary chalcopyrite materials, said single source precursor being a liquid at room temperature and being effective to yield a ternary chalcopyrite material upon heating or pyrolysis thereof.
- 39. (original) A single source precursor according to claim 38, said single source precursor being effective to yield a I-III-VI₂ ternary chalcopyrite material upon heating or pyrolysis of said single source precursor at a temperature less than about 500°C.
- 40. (currently amended) A method of depositing ternary chalcopyrite materials comprising the steps of:
- a) providing a first single source precursor for said ternary chalcopyrite material, said first single source precursor having the empirical formula $[\{L\}_nM'(ER)_x(X)_y(R)_zM'']$, wherein x is 1-4, x+y+z=4, n is greater than or equal to 1, L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom, M' is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl,

vinyl, perfluoro alkyl, perfluoro aryl, silane, and carbamato groups, said single source precursor excluding

$$\begin{split} & \left[\{ P(C_6H_5)_3 \}_2 Cu(S-C_2H_5)_2 In(S-C_2H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Cu(SC\{O\}C_5H_6)_2 In(SC\{O\}C_5H_6)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Cu(SC\{O\}C_5H_6)_2 Ga(SC\{O\}C_5H_6)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}C_5H_6)_2 In(SC\{O\}C_5H_6)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}C_5H_6)_2 Ga(SC\{O\}C_5H_6)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right] \\ & \left[\{ P(C_6H_5)_3 \}_2 Cu(S-C_2H_5)_2 In(S-C_2H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Cu(SC\{O\}C_6H_5)_2 In(SC\{O\}C_6H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Cu(SC\{O\}C_6H_5)_2 In(SC\{O\}C_6H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}C_6H_5)_2 In(SC\{O\}C_6H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}C_6H_5)_2 In(SC\{O\}C_6H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}C_6H_5)_2 In(SC\{O\}C_6H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}C_6H_5)_2 In(SC\{O\}C_6H_5)_2 \right], \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right], \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & and \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right]; \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right] \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right] \\ & \left[\{ P(C_6H_5)_3 \}_2 Ag(SC\{O\}CH_3)_2 In(SC\{O\}CH_3)_2 \right] \\ & \left[\{ P(C_6H_5$$

- b) depositing the single source precursor on a substrate using a spray CVD technique.
- 41. (original) A method according to claim 40, said single source precursor having a structural formula selected from the group consisting of



42. (original) A method according to claim 40, said single source precursor having a structural formula selected from the group consisting of

$$L \longrightarrow M$$

$$\downarrow E$$

$$\downarrow R$$

43. (original) A method according to claim 40, said single source precursor having a structural formula selected from the group consisting of

$$\begin{array}{c|c} & & & & \\ & &$$

44. (original) A method according to claim 40, said single source precursor having a structural formula selected from the group consisting of

$$L \longrightarrow M! \longrightarrow R$$
 and
$$L \longrightarrow M! \longrightarrow R$$

$$R$$

45. (original) A method according to claim 40, said single source precursor having three E-R groups.

- 46. (original) A method according to claim 40, comprising the steps of providing a second single source precursor, and applying said first and second single source precursors on said substrate via said spray CVD technique.
- 47. (currently amended) A method of making a single source precursor for the deposition of ternary chalcopyrite materials comprising the step of carrying out the following reaction:

$$4M_{ALK}ER + M'X_3 + M''X + nL \longrightarrow [\{L\}_nM'(ER)_2M''(ER)_2]$$
 wherein

M_{ALK} is an alkali metal element,

E is a Group VI-A element,

R is selected from the group consisting of alkyl, aryl, vinyl, perfluoro alkyl, perfluoro aryl, silane and carbamato groups,

M' is a Group [[III-A]] I-B element,

M" is a Group [[I-B]] III-A element,

X is a Group VII-A element, and

n is greater than or equal to 1.

- 48. (original) A method according to claim 47, wherein said single source precursor is made in a single step consisting essentially of said reaction.
- 49. (original) A method according to claim 47, wherein the ionic complex $[L_{(n)}M"(CH_3CN)_{(4-n)}]^+$ is formed *in situ* as said reaction proceeds.
- 50. (original) A method according to claim 47, said reaction being carried out under anaerobic conditions.
- 51. (original) A method according to claim 47, said reaction being carried out under non-anaerobic conditions.
 - 52. (original) A method of making a quantum dot comprising the steps of:
 - a) providing a single source precursor for a ternary chalcopyrite material; and
- b) pyrolyzing said single source precursor to yield a quantum dot made of ternary chalcopyrite material having dimensions less than 100 nanometers.

- 53. (original) A method according to claim 52, said quantum dot made of a ternary I-III-VI₂ chalcopyrite material.
- 54. (original) A method according to claim 52, said quantum dot made of a ternary I-III₅-VI₈ chalcopyrite material.
- 55. (original) A method according to claim 52, said pyrolyzing step being carried out at a temperature less than about 500°C.
- 56. (original) A method according to claim 52, said single source precursor having the empirical formula $[\{L\}_nM'(ER)_x(X)_y(R)_zM'']$, wherein x is 1-4, x+y+z=4, n is greater than or equal to 1, L is a Lewis base that is coordinated to M' via a dative bond, M' is a Group I-B atom, M' is a Group III-A atom, E is a Group VI-A atom, X is a Group VII-A atom, and each R is individually selected from the group consisting of alkyl, aryl, vinyl, perfluoro alkyl, perfluoro aryl, silane, and carbamato groups.
- 57. (original) A single source precursor according to claim 2, said single source precursor being effective to yield a I-III₅-VI₈ ternary chalcopyrite material upon heating or pyrolysis of said single source precursor.
- 58. (original) A single source precursor according to claim 30, said single source precursor being effective to yield a ternary chalcopyrite material having a band gap of 0.5-3.5 eV between a conduction band and a valence band thereof, said ternary chalcopyrite material being (Cu:Ag:Au)₁(Al:In:Ga)₁(S:Se:Te)₂.
- 59. (new) A single source precursor according to claim 2, said Group VI-A atom being selected from the group consisting of S, Se and Te.
- 60. (new) A single source precursor according to claim 16, said Group VI-A atom being selected from the group consisting of S, Se and Te.